

Data Path : V:\HPCHEM1\BNA M\DATA\BM070916\
 Data File : BM006356.D
 Acq On : 08 Jul 2016 18:42
 Operator : UM/SJ
 Sample : H3861-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BDJ00

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM070216.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.828	36	41	53	rVB	155322	312730	2.10%	0.382%
2	6.816	712	719	729	rBV	505040	823075	5.53%	1.006%
3	6.981	740	747	760	rVB	362901	586439	3.94%	0.717%
4	7.175	773	780	791	rVB	500889	810407	5.44%	0.990%
5	7.639	851	859	867	rBV	735021	1187175	7.97%	1.451%
6	8.345	973	979	995	rVB	570758	985619	6.62%	1.204%
7	8.792	1048	1055	1065	rBV	512033	838564	5.63%	1.025%
8	9.510	1169	1177	1187	rBV	406467	716135	4.81%	0.875%
9	10.045	1260	1268	1281	rBV	598677	1067269	7.17%	1.304%
10	10.422	1321	1332	1345	rBV	1012144	1837926	12.35%	2.246%
11	10.563	1348	1356	1367	rBV	568165	1080159	7.26%	1.320%
12	13.698	1883	1889	1894	rBV	1119694	1691108	11.36%	2.066%
13	13.745	1894	1897	1907	rVB	687323	1047329	7.03%	1.280%
14	13.980	1929	1937	1947	rBV	1242503	2033331	13.66%	2.484%
15	14.292	1977	1990	1999	rBV2	1431135	2441867	16.40%	2.984%
16	14.510	2021	2027	2039	rBV	285371	545464	3.66%	0.666%
17	15.286	2152	2159	2170	rBV	1734681	2522400	16.94%	3.082%
18	15.415	2176	2181	2194	rVB	289774	431864	2.90%	0.528%
19	17.039	2451	2457	2467	rBV	1899861	2861217	19.22%	3.496%
20	17.139	2467	2474	2486	rVB2	1776605	2672239	17.95%	3.265%
21	19.439	2859	2865	2874	rBV2	1986955	2919572	19.61%	3.567%
22	20.374	3020	3024	3032	rVB	430134	537846	3.61%	0.657%
23	20.491	3039	3044	3047	rBV	608656	748851	5.03%	0.915%
24	20.556	3047	3055	3061	rVB2	2668267	4317396	29.00%	5.275%
25	20.721	3078	3083	3089	rBV	1133366	1542002	10.36%	1.884%
26	20.786	3089	3094	3099	rVV	3004202	4065244	27.31%	4.967%
27	20.839	3099	3103	3108	rVV2	2432046	3653863	24.54%	4.465%
28	20.897	3108	3113	3118	rVB4	839144	1670789	11.22%	2.041%
29	21.091	3140	3146	3151	rBV4	5433804	9777780	65.68%	11.947%
30	21.150	3151	3156	3161	rVV	10179307	14887972	100.00%	18.191%
31	21.197	3161	3164	3167	rVV	1103522	1326112	8.91%	1.620%
32	21.238	3167	3171	3179	rVV	1799337	2412434	16.20%	2.948%
33	21.333	3179	3187	3193	rVV	1007301	1316903	8.85%	1.609%
34	21.450	3203	3207	3212	rVV	661595	867165	5.82%	1.060%

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35	23.303	3516	3522	3537	rVB2	1074798	2220340	14.91%	2.713%
36	23.444	3539	3546	3560	rVB	1120323	2234895	15.01%	2.731%
37	23.956	3629	3633	3643	rVB2	140301	257153	1.73%	0.314%
38	24.562	3731	3736	3743	rVB3	157761	311982	2.10%	0.381%
39	24.685	3753	3757	3769	rVB2	125064	280960	1.89%	0.343%

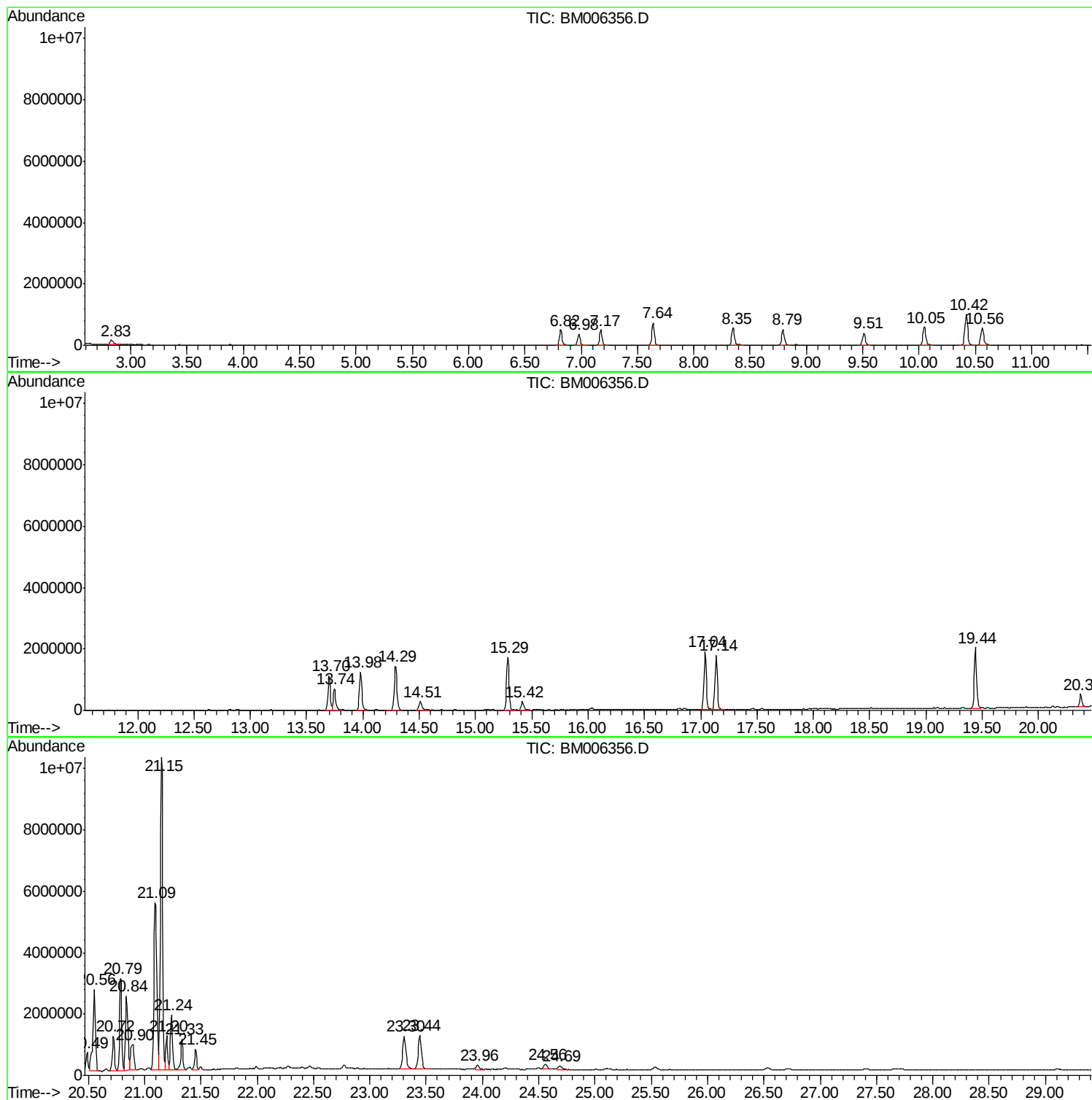
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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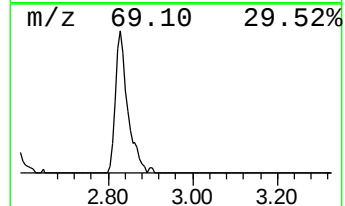
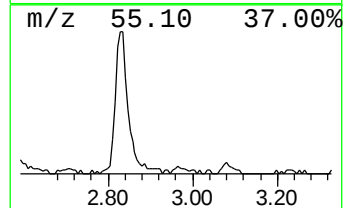
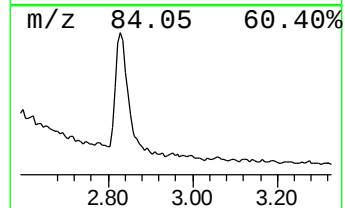
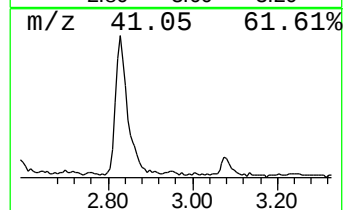
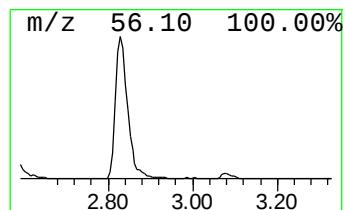
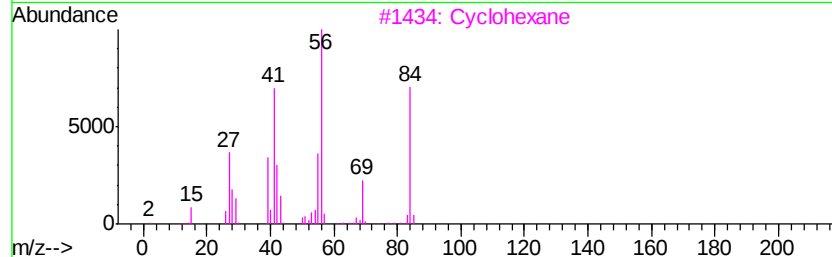
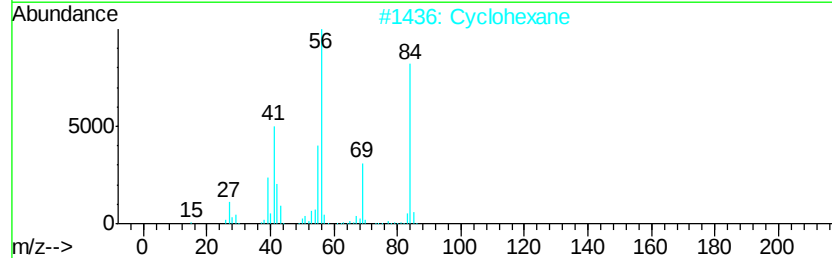
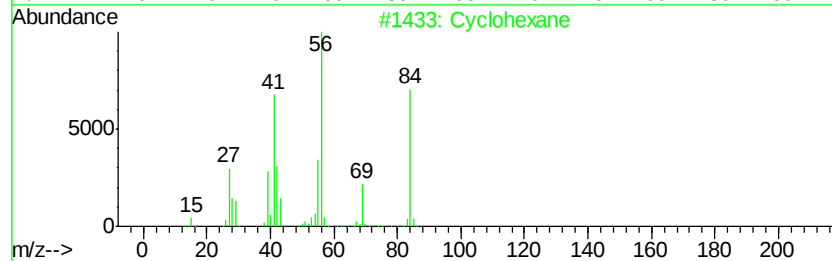
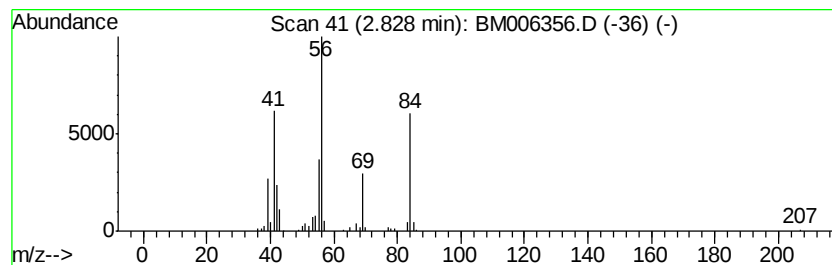
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM070216.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Cyclic2.83 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.83	5.27 ng/ul	312730	1,4-Dichlorobenzene-d4	7.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane	84	C6H12	000110-82-7	91
2		Cyclohexane	84	C6H12	000110-82-7	91
3		Cyclohexane	84	C6H12	000110-82-7	91
4		Cyclohexane	84	C6H12	000110-82-7	91
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	59



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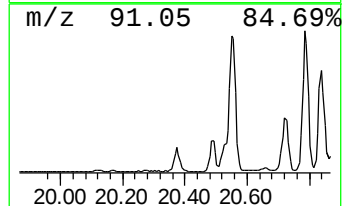
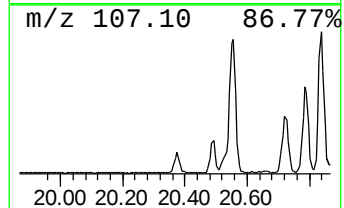
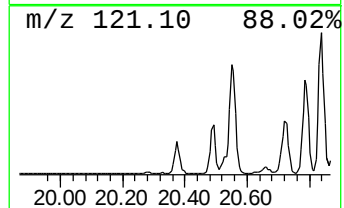
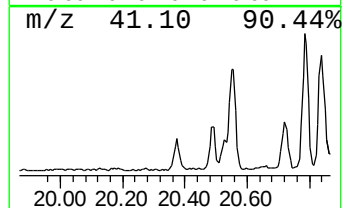
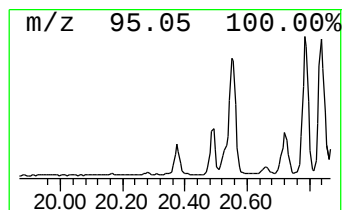
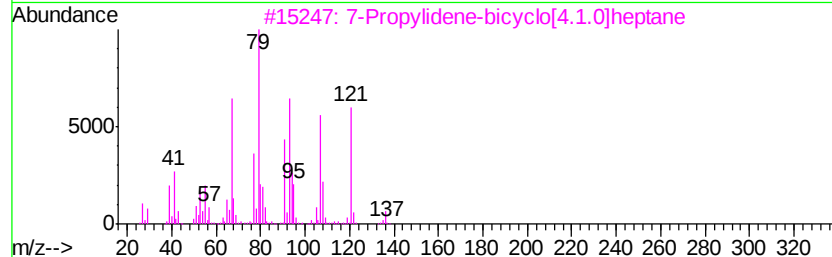
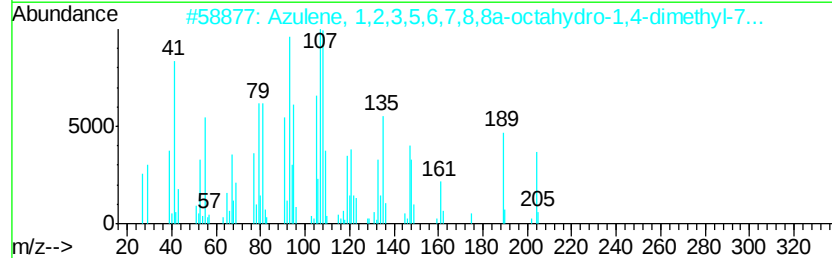
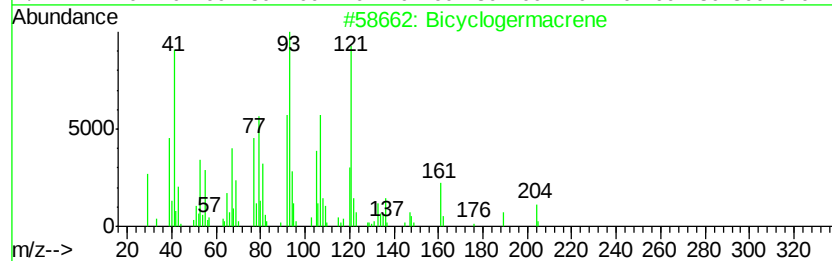
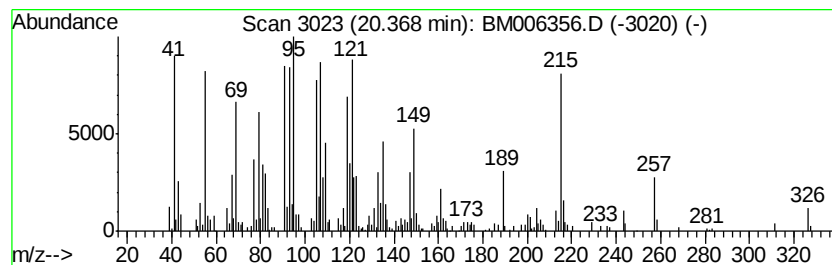
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Peak Number 2 Bicyclogermacrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.37	4.46 ng/ul	537846	Chrysene-d12	21.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Bicyclogermacrene	204 C15H24	067650-90-2 52
2		Azulene, 1,2,3,5,6,7,8,8a-octahy...	204 C15H24	003691-11-0 43
3		7-Propylidene-bicyclo[4.1.0]heptane	136 C10H16	082253-09-6 38
4		1H-Cycloprop[el]azulene, 1a,2,3,5...	204 C15H24	021747-46-6 30
5		1-Cycloheptene, 1,4-dimethyl-3-(...	204 C15H24	1000159-38-6 25



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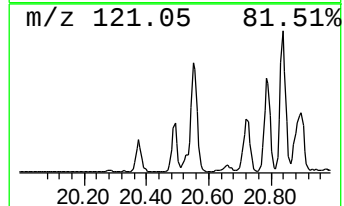
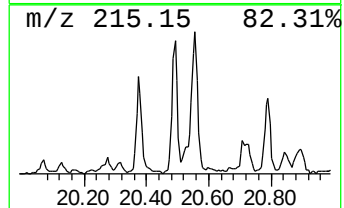
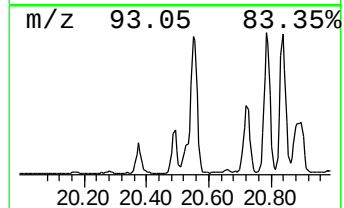
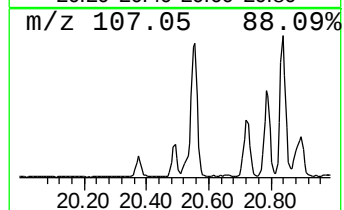
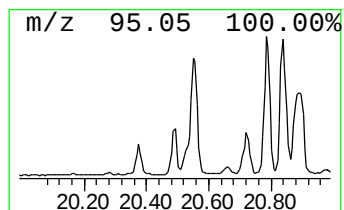
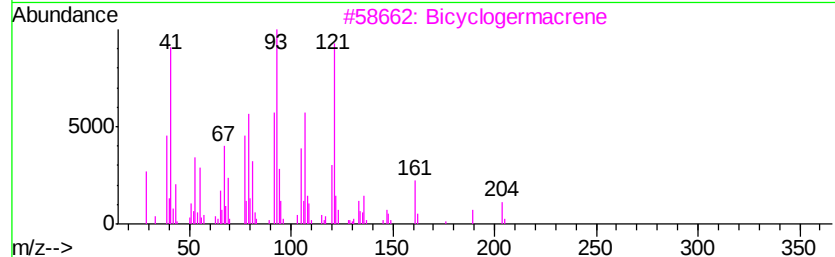
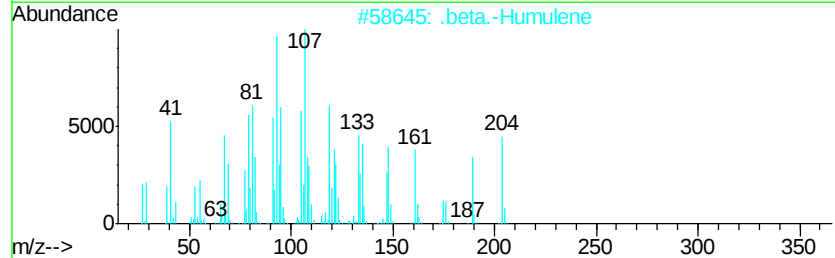
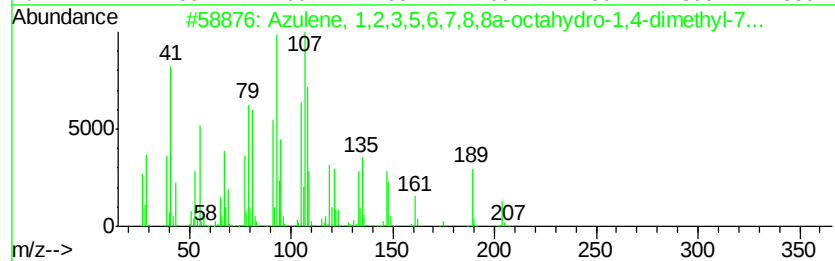
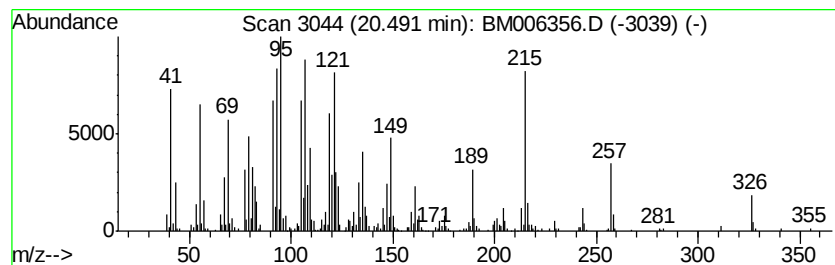
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Peak Number 3 Azulene, 1,2,3,5,6,7,8,8a-o... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.49	6.21 ng/ul	748851	Chrysene-d12	21.24		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Azulene, 1,2,3,5,6,7,8,8a-octahy...	204	C15H24	003691-11-0	50
2		.beta.-Humulene	204	C15H24	000116-04-1	42
3		Bicyclogermacrene	204	C15H24	067650-90-2	38
4		.alpha.-Carvophyllene	204	C15H24	006753-98-6	30
5		1H-Cycloprop[e]azulene, 1a,2,3,5...	204	C15H24	021747-46-6	27



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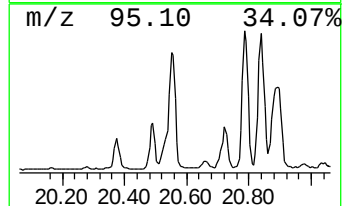
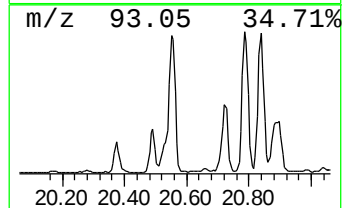
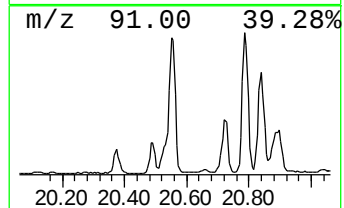
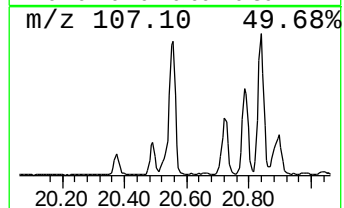
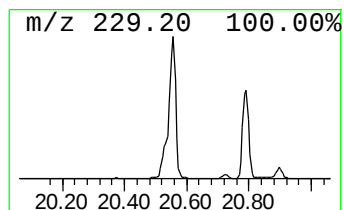
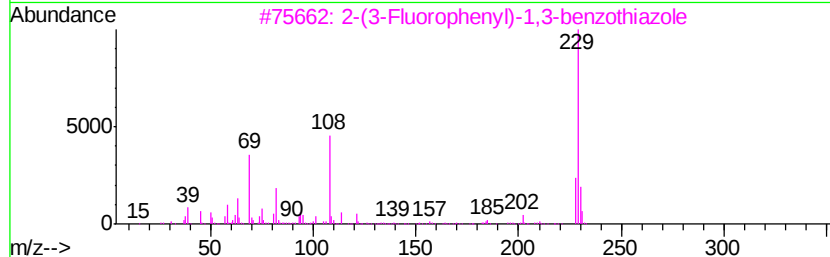
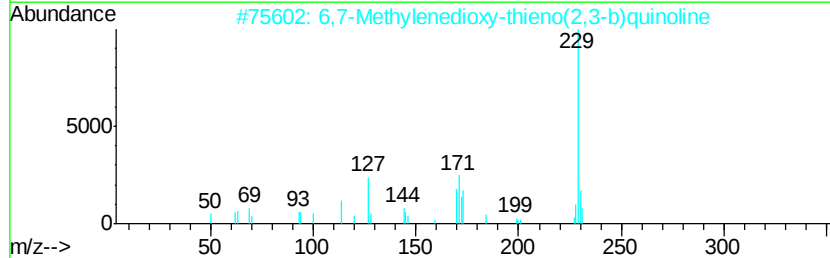
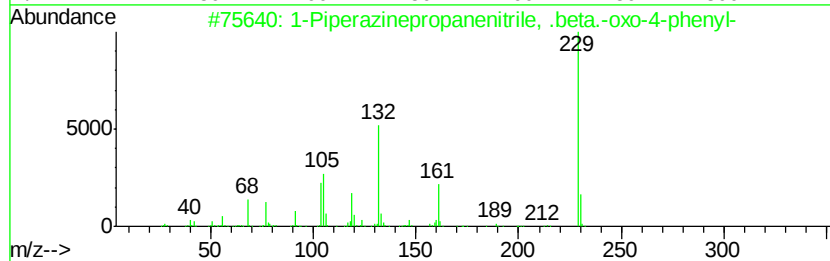
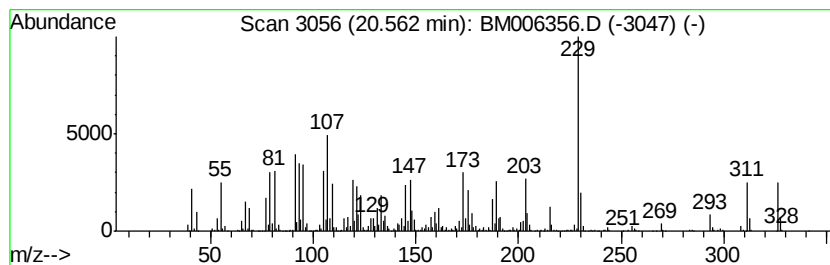
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Peak Number 4 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.56	35.79 ng/ul	4317400	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Piperazinepropanenitrile, .bet...	229	C13H15N3O	014761-40-1	25
2		6,7-Methylenedioxy-thieno(2,3-b)...	229	C12H7N02S	062638-09-9	25
3		2-(3-Fluorophenyl)-1,3-benzothia...	229	C13H8FNS	1000298-18-6	22
4		Benzoic acid, 4-(4-propylcyclohe...	381	C25H35NO2	1000204-13-6	22
5		10H-Phenothiazine, 1-methoxy-	229	C13H11NOS	001576-70-1	22



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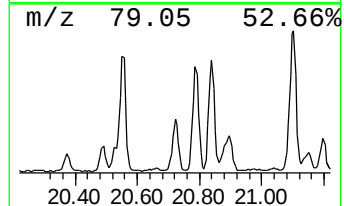
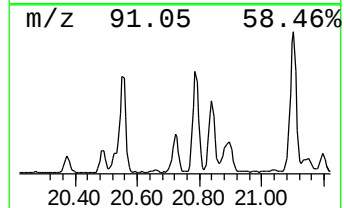
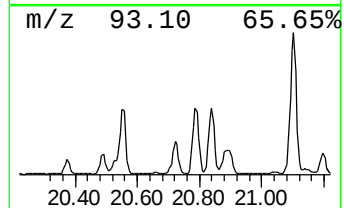
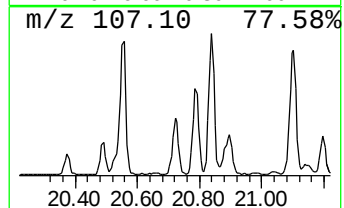
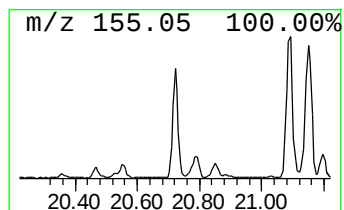
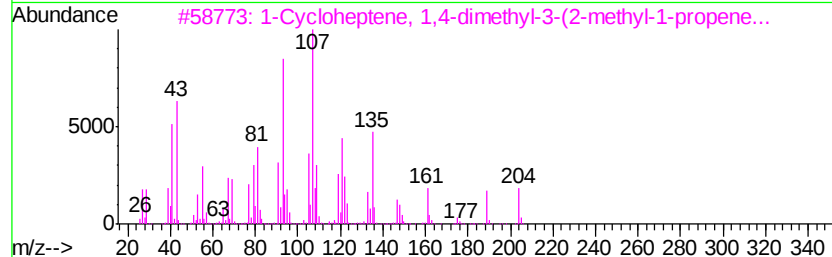
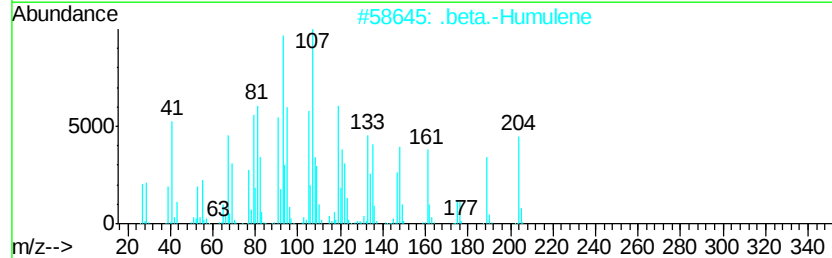
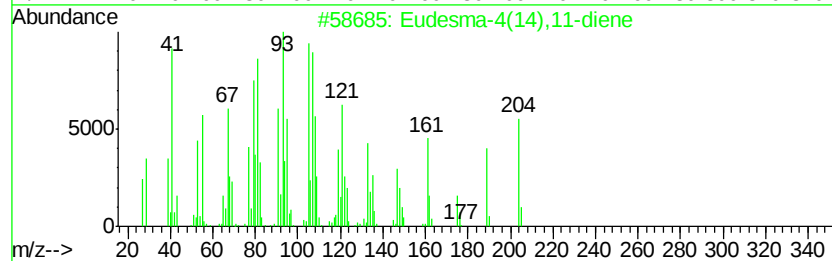
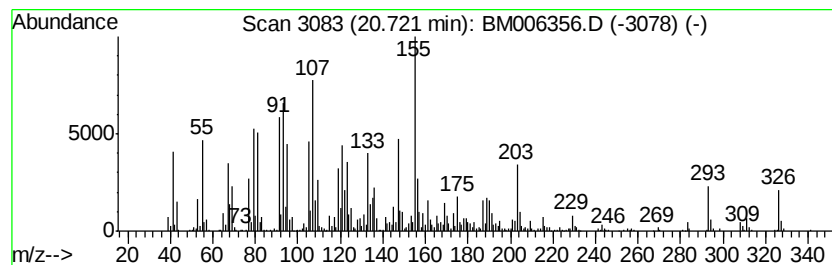
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Peak Number 5 Eudesma-4(14),11-diene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.72	12.78 ng/ul	1542000	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eudesma-4(14),11-diene	204	C15H24	1000152-04-3	59
2		.beta.-Humulene	204	C15H24	000116-04-1	38
3		1-Cycloheptene, 1,4-dimethyl-3-(...	204	C15H24	1000159-38-6	25
4		1,3-Dimethyl-5-(propen-1-yl)adam...	204	C15H24	057040-48-9	25
5		2-(6-Methylpyridin-2-ylmethyl)cy...	205	C13H19NO	128231-64-1	25



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Data File : BM006356.D
Acq On : 08 Jul 2016 18:42
Operator : UM/SJ
Sample : H3861-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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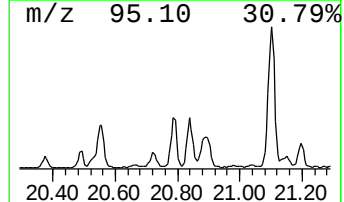
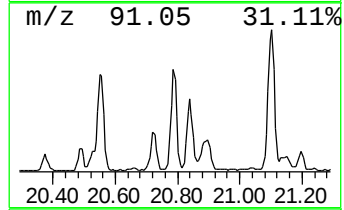
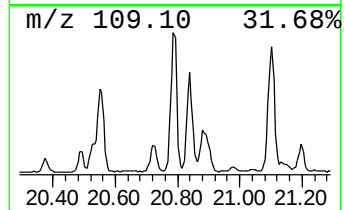
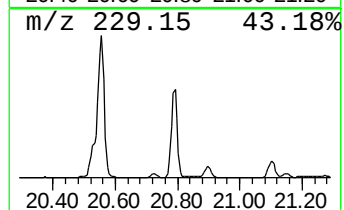
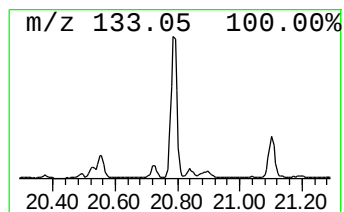
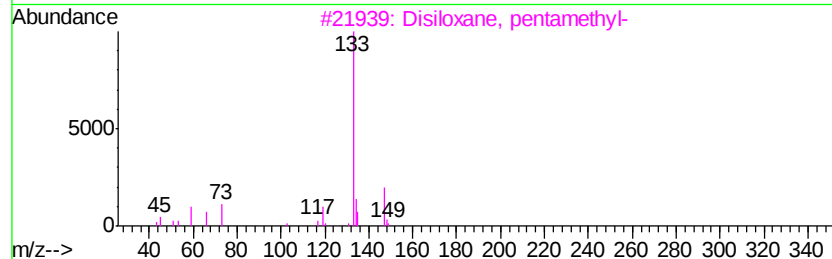
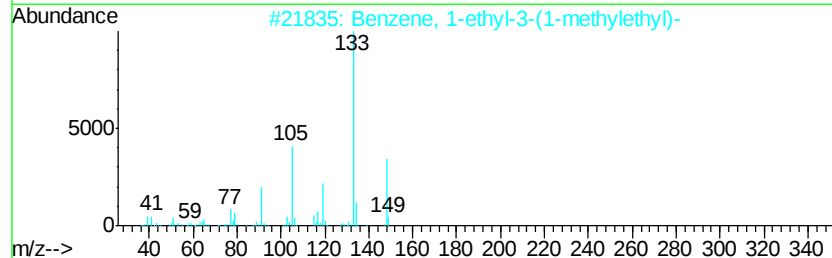
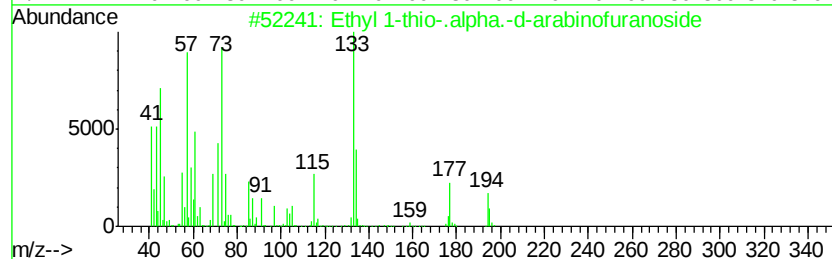
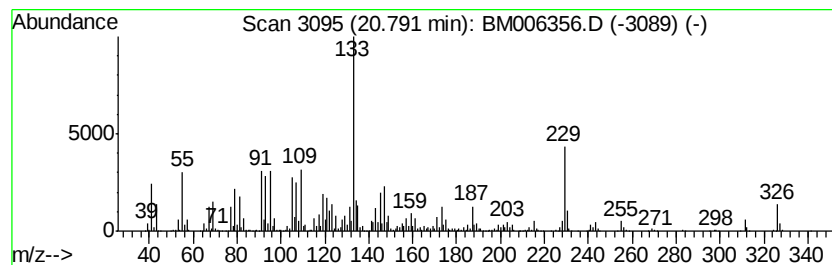
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Ethyl 1-thio-.alpha.-d-arab... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.79	33.70 ng/ul	4065240	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethyl 1-thio-.alpha.-d-arabinoфу...	194	C7H14O4S	1000129-83-7	60
2		Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	38
3		Disiloxane, pentamethyl-	148	C5H16OSi2	001438-82-0	27
4		2-{2-[1-(2-Hydroxyethyl)-5-oxo-3...}	417	C20H23N3O5S	1000224-26-4	25
5		Benzeneacetonitrile, 4-hydroxy-	133	C8H7NO	014191-95-8	22



Data Path : V:\HPCHEM1\BNA M\DATA\BM070916\
Data File : BM006356.D
Acq On : 08 Jul 2016 18:42
Operator : UM/SJ
Sample : H3861-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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ClientSampleId :
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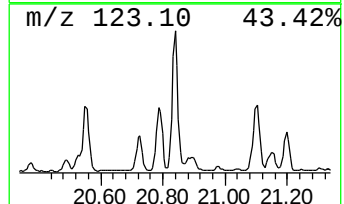
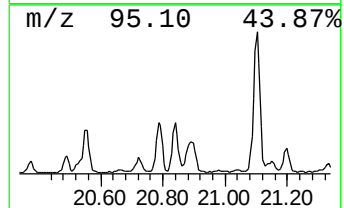
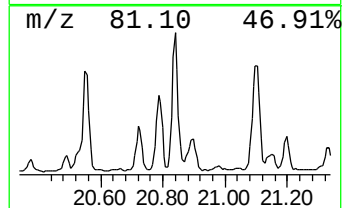
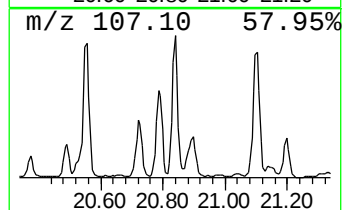
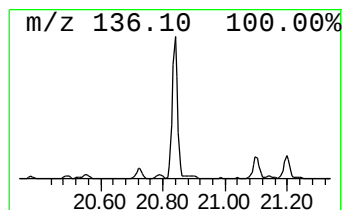
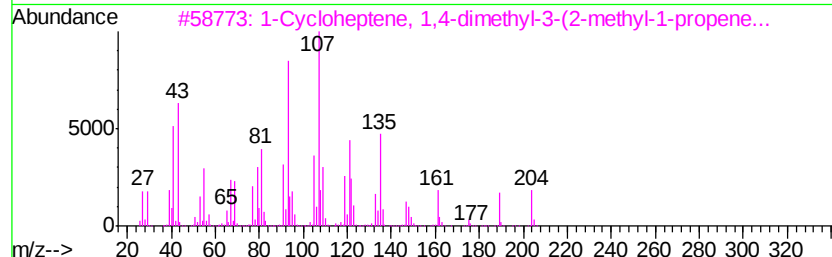
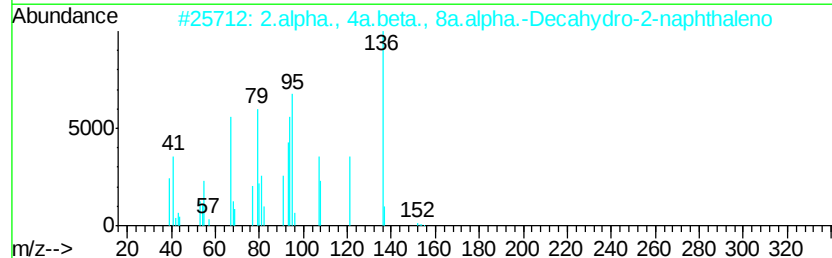
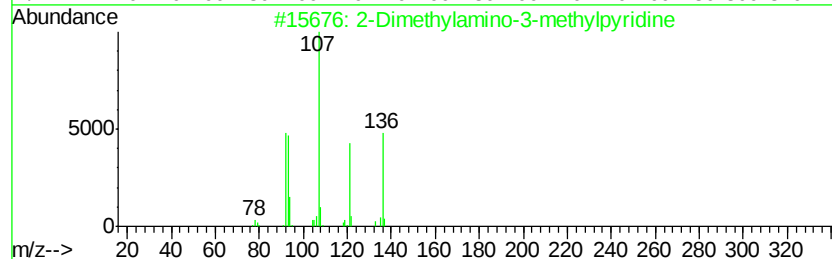
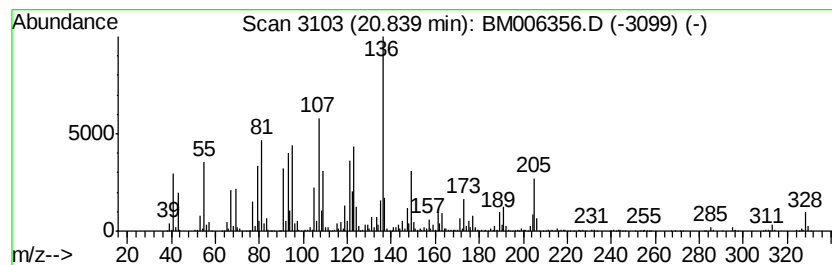
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 2-Dimethylamino-3-methylpyr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.84	30.29 ng/ul	3653860	Chrysene-d12	21.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Dimethylamino-3-methylpyridine	136	C8H12N2	061713-46-0	50
2		2.alpha., 4a.beta., 8a.alpha.-De...	154	C10H18O	005779-35-1	41
3		1-Cycloheptene, 1,4-dimethyl-3-(...	204	C15H24	1000159-38-6	25
4		Benzenemethanol, 3,5-dimethyl-	136	C9H12O	027129-87-9	25
5		2-Hydroxy-5-methylbenzaldehyde	136	C8H8O2	000613-84-3	22



Data Path : V:\HPCHEM1\BNA M\DATA\BM070916\
Data File : BM006356.D
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Sample : H3861-01
Misc :
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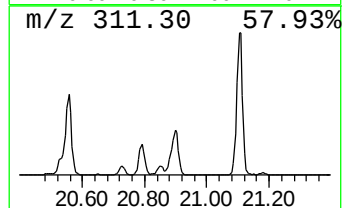
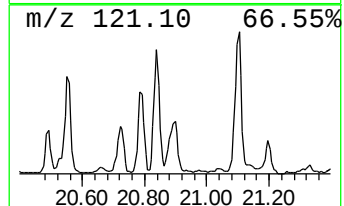
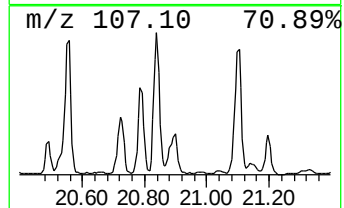
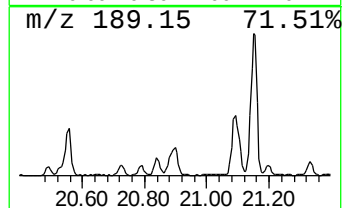
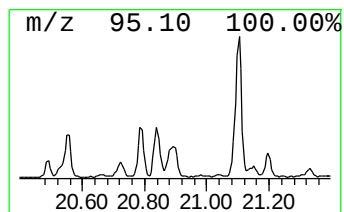
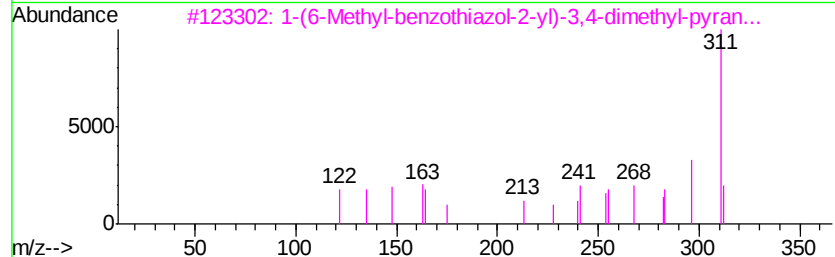
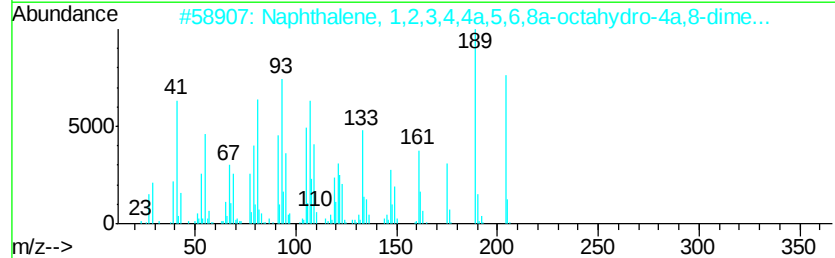
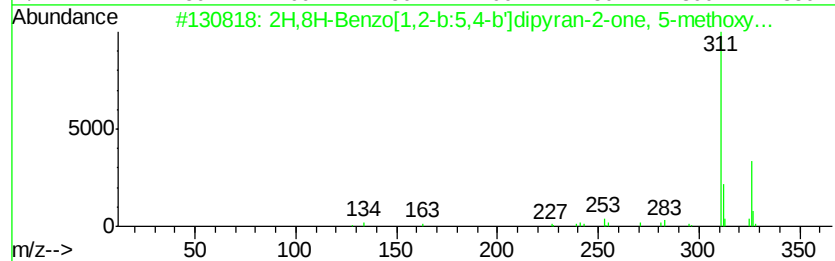
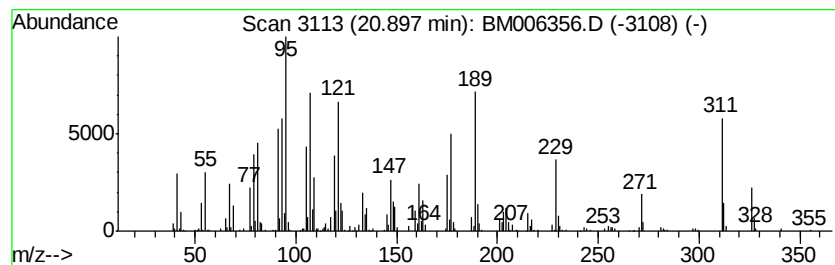
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.90	13.85 ng/ul	1670790	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H,8H-Benzo[1,2-b:5,4-b']dipyran...	326	C20H22O4	055334-39-9	25
2		Naphthalene, 1,2,3,4,4a,5,6,8a-o...	204	C15H24	000473-13-2	25
3		1-(6-Methyl-benzothiazol-2-yl)-3...	311	C16H13N3O2S	106531-84-4	11
4		Azulene, 1,2,3,5,6,7,8,8a-octahy...	204	C15H24	003691-11-0	11
5		1-Cycloheptene, 1,4-dimethyl-3-(...	204	C15H24	1000159-38-6	11



Data Path : V:\HPCHEM1\BNA M\DATA\BM070916\
Data File : BM006356.D
Acq On : 08 Jul 2016 18:42
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Sample : H3861-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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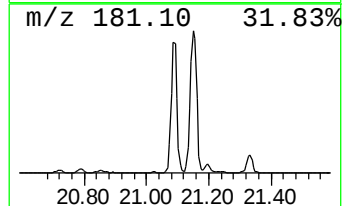
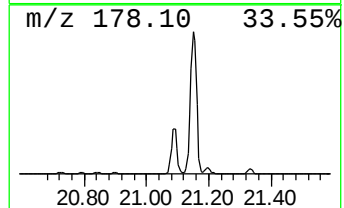
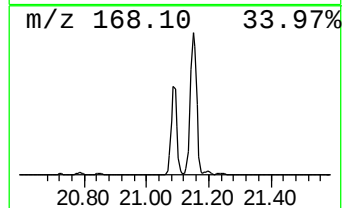
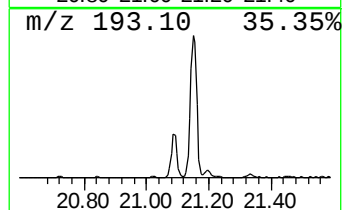
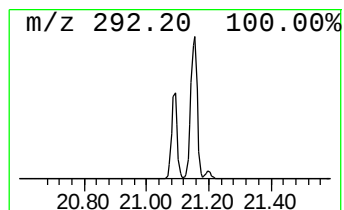
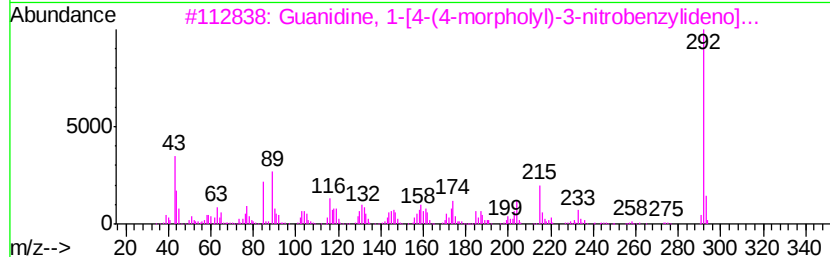
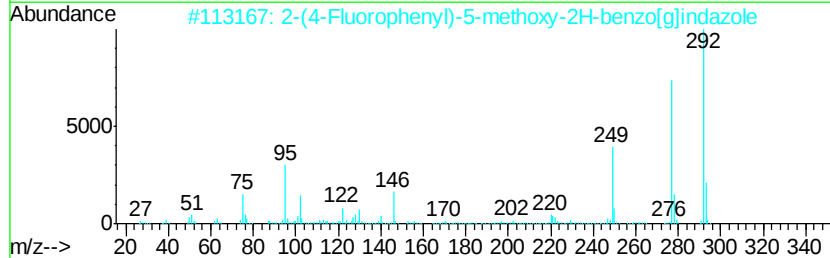
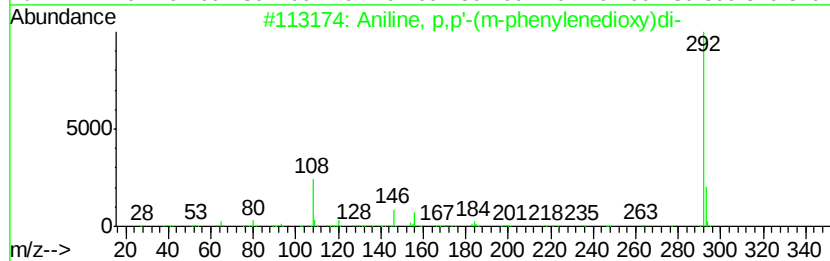
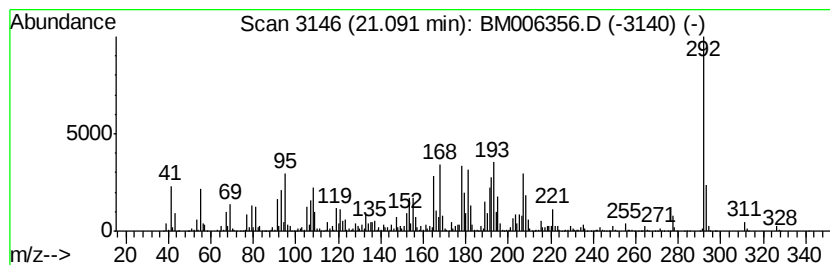
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.09	81.06 ng/ul	9777780	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Aniline, p,p'-(m-phenylenedioxy)di-	292	C18H16N2O2	002479-46-1	30
2		2-(4-Fluorophenyl)-5-methoxy-2H-...	292	C18H13FN2O	1000194-51-1	30
3		Guanidine, 1-[4-(4-morpholyl)-3-...	292	C12H16N6O3	296244-85-4	25
4		1,1'-Biphenyl, 4,4-bis(1-pyrroli...	292	C20H24N2	1000193-54-6	25
5		Benzoic acid, 4-[2-(2-hydroxy-1-...	292	C17H12N2O3	1000197-03-1	22



Data Path : V:\HPCHEM1\BNA M\DATA\BM070916\
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Operator : UM/SJ
Sample : H3861-01
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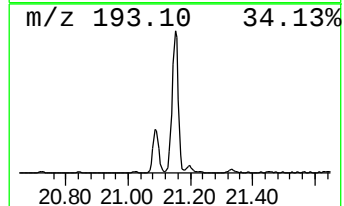
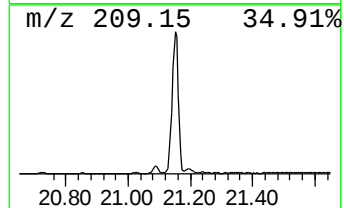
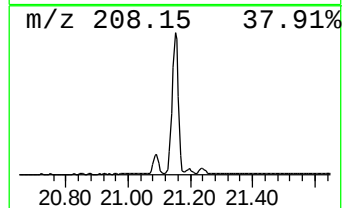
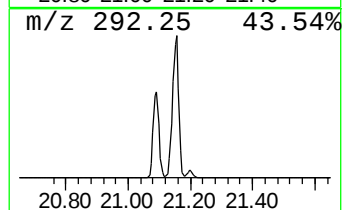
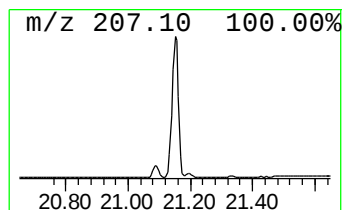
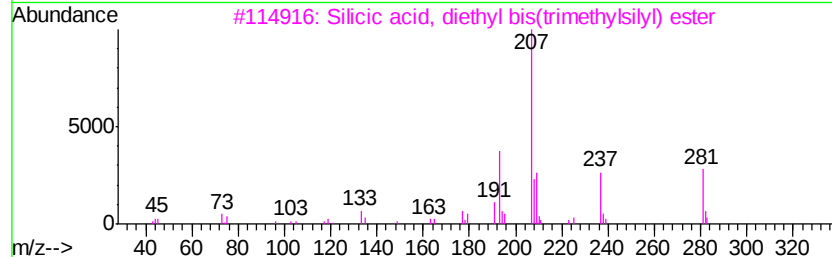
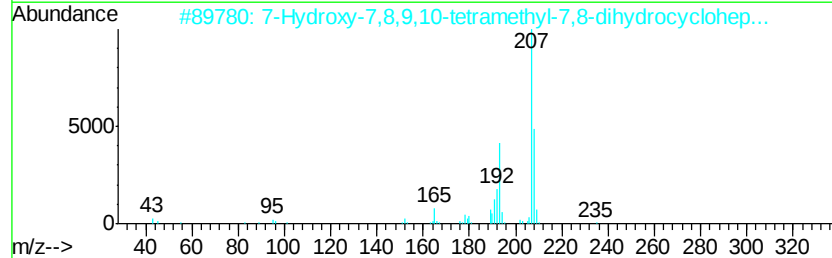
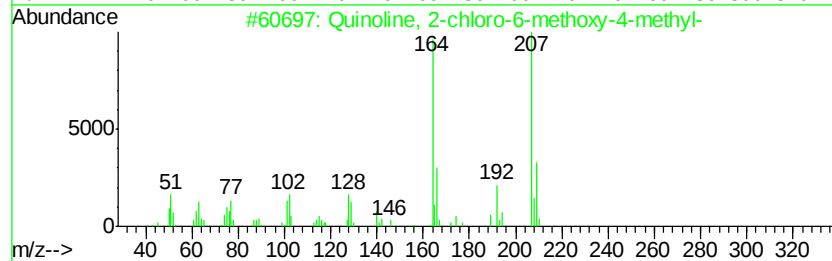
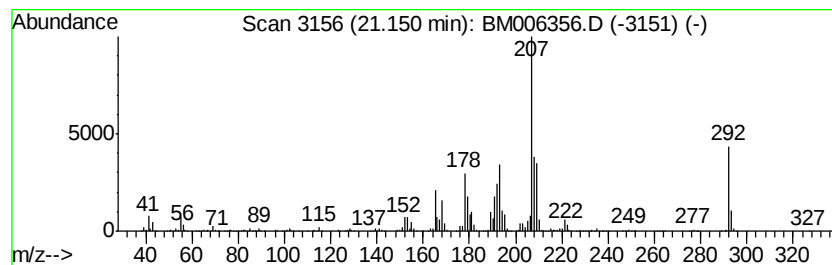
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM070216.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown-04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.15	123.43 ng/ul	14888000	Chrysene-d12	21.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Quinoline, 2-chloro-6-methoxy-4-...	207 C11H10ClNO	006340-55-2	43
2	7-Hydroxy-7,8,9,10-tetramethyl-7...	252 C18H20O	1000110-34-9	43
3	Silicic acid, diethyl bis(trimet...	296 C10H28O4Si3	003555-45-1	37
4	7-Chloro-4-methoxy-3-methylquino...	207 C11H10ClNO	1000213-52-2	37
5	1-Methyl-3-phenylindole	207 C15H13N	030020-98-5	32



Data Path : V:\HPCHEM1\BNA M\DATA\BM070916\
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Misc :
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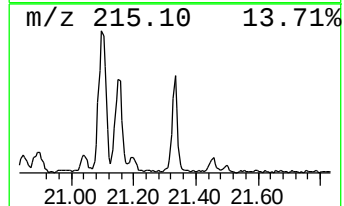
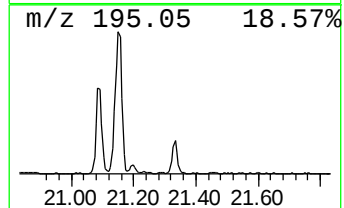
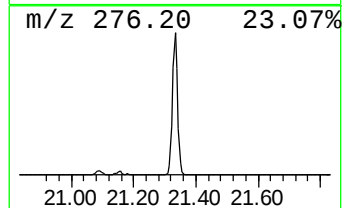
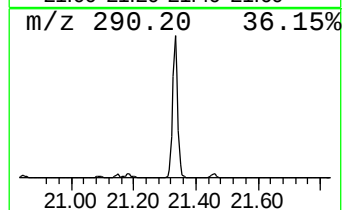
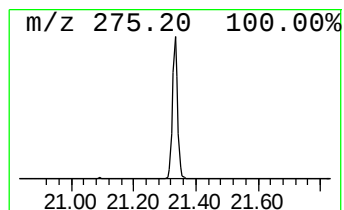
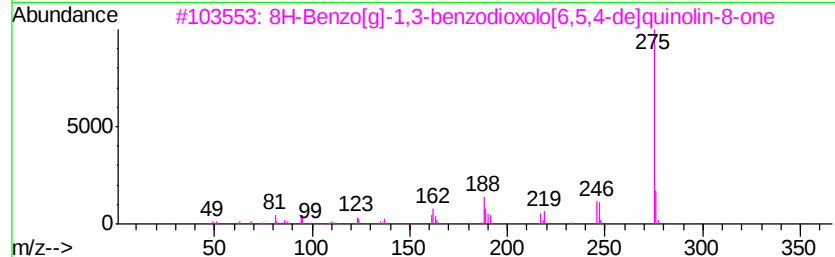
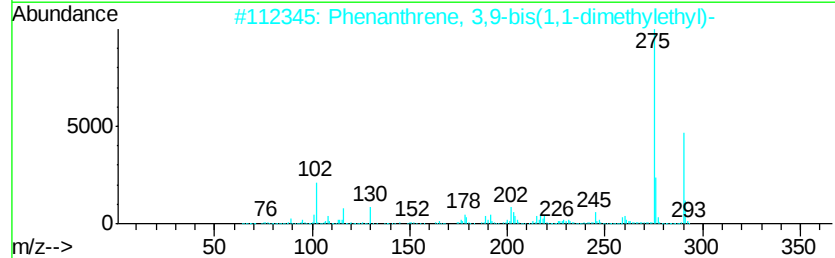
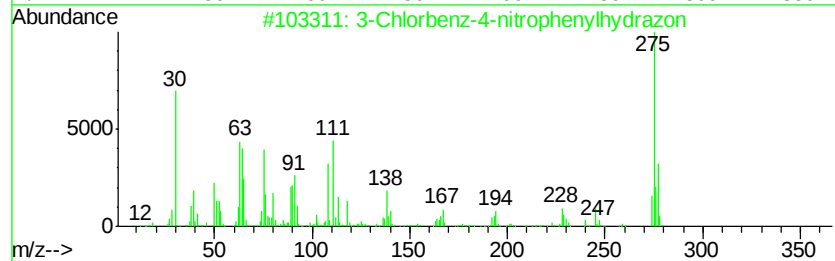
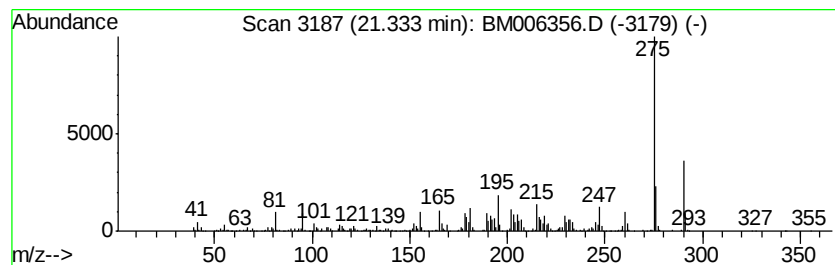
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 3-Chlorbenz-4-nitrophenylhy... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.33	10.92 ng/ul	1316900	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chlorbenz-4-nitrophenylhydrazon	275	C13H10ClN3O2	005908-26-9	91
2		Phenanthrene, 3,9-bis(1,1-dimeth...	290	C22H26	055125-03-6	87
3		8H-Benzo[<i>a</i>]-1,3-benzodioxolo[6,5...	275	C17H9NO3	000475-75-2	64
4		8H-Benzo[<i>a</i>]-1,3-benzodioxolo[6,5...	275	C17H9NO3	000475-75-2	62
5		4,4'-(Alpha-methylbenzylidene)di...	290	C20H18O2	001571-75-1	59



Data Path : V:\HPCHEM1\BNA M\DATA\BM070916\
Data File : BM006356.D
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ALS Vial : 4 Sample Multiplier: 1

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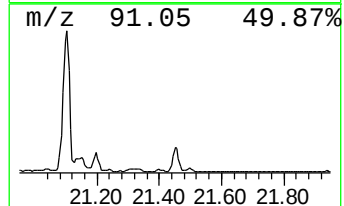
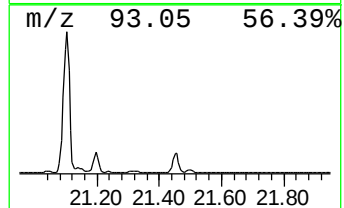
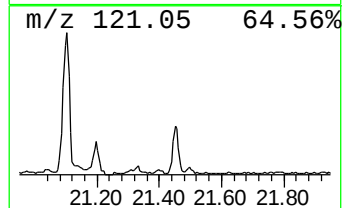
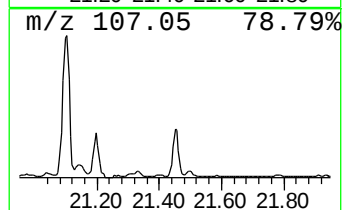
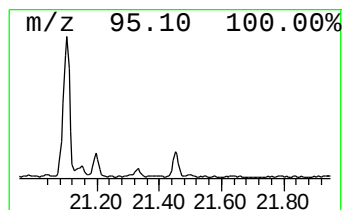
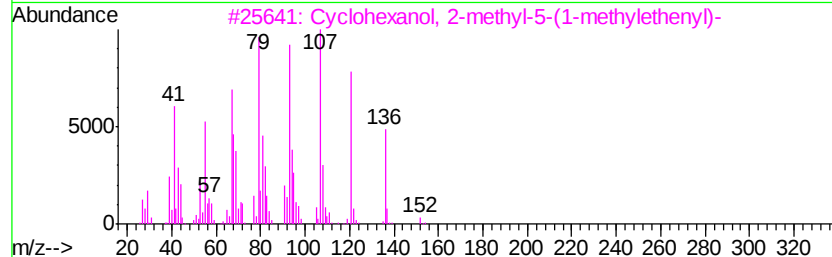
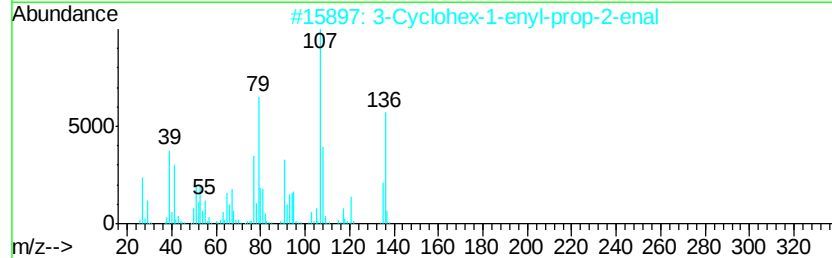
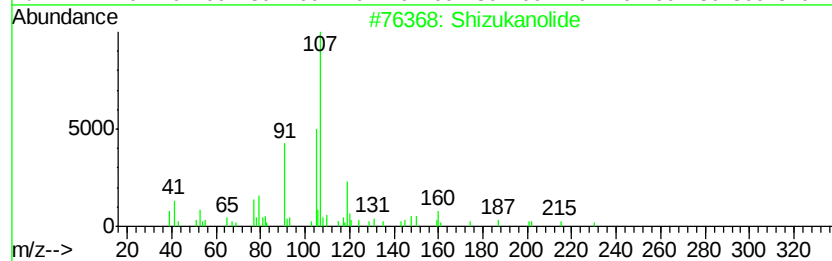
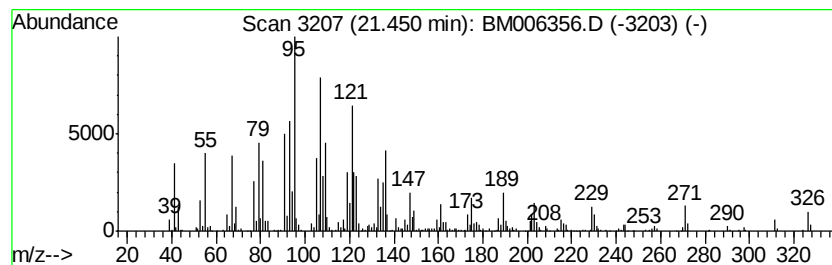
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown-05 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.45	7.19 ng/ul	867165	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Shizukanolide	230	C15H18O2	070578-36-8	44
2		3-Cyclohex-1-enyl-prop-2-enal	136	C9H12O	1000193-70-2	38
3		Cyclohexanol, 2-methyl-5-(1-meth...	154	C10H18O	000619-01-2	38
4		Cyclohexane, 1,2-dimethyl-3,5-bi...	192	C14H24	074806-55-6	35
5		2-(6-Methylpyridin-2-ylmethyl)cy...	205	C13H19NO	128231-64-1	25



Data Path : V:\HPCHEM1\BNA M\DATA\BM070916\
Data File : BM006356.D
Acq On : 08 Jul 2016 18:42
Operator : UM/SJ
Sample : H3861-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BDJ00

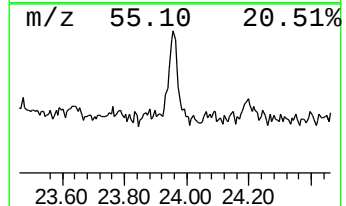
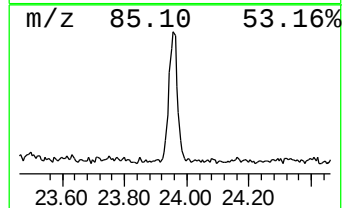
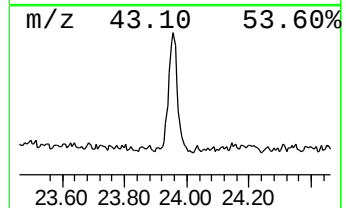
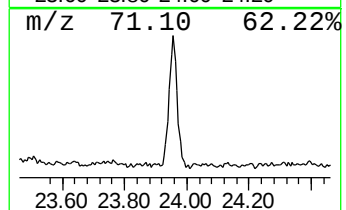
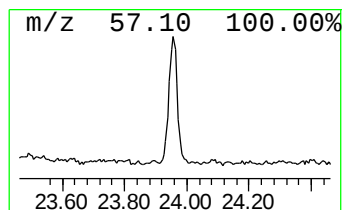
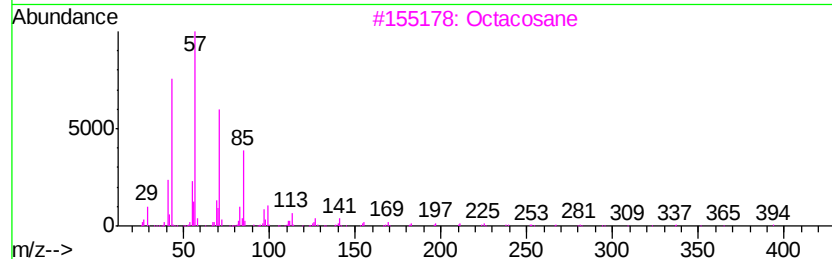
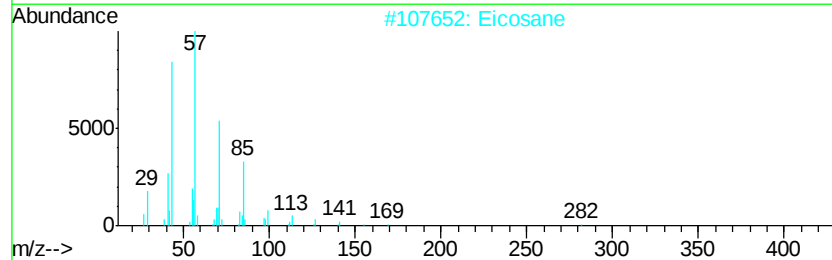
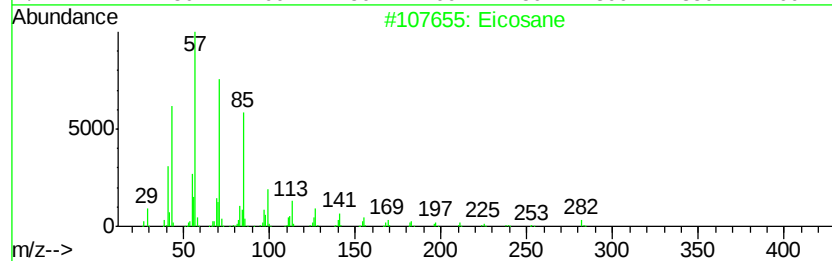
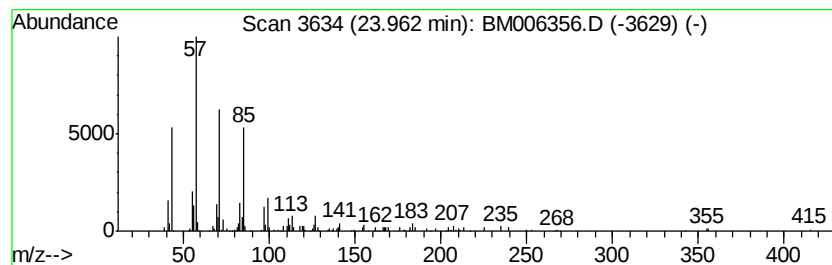
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.96	2.30 ng/ul	257153	Perylene-d12	23.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	93
2	Eicosane			282	C20H42	000112-95-8	92
3	Octacosane			394	C28H58	000630-02-4	90
4	Heptadecane			240	C17H36	000629-78-7	90
5	Octacosane			394	C28H58	000630-02-4	90



Data Path : V:\HPCHEM1\BNA M\DATA\BM070916\
Data File : BM006356.D
Acq On : 08 Jul 2016 18:42
Operator : UM/SJ
Sample : H3861-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BDJ00

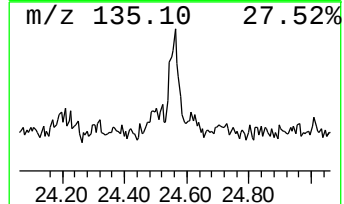
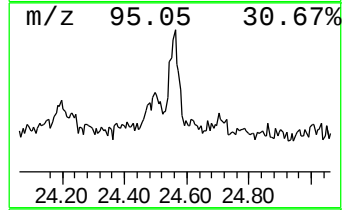
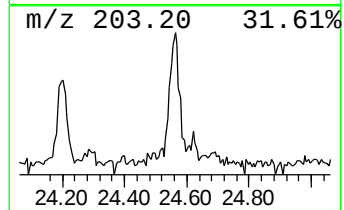
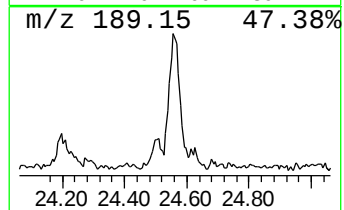
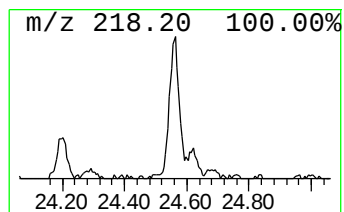
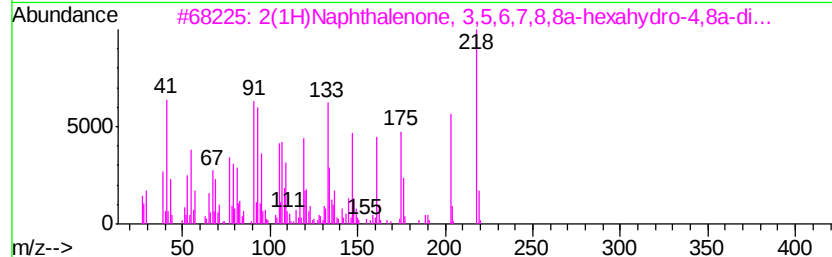
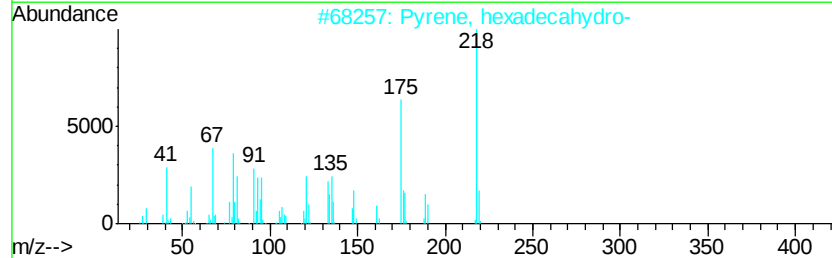
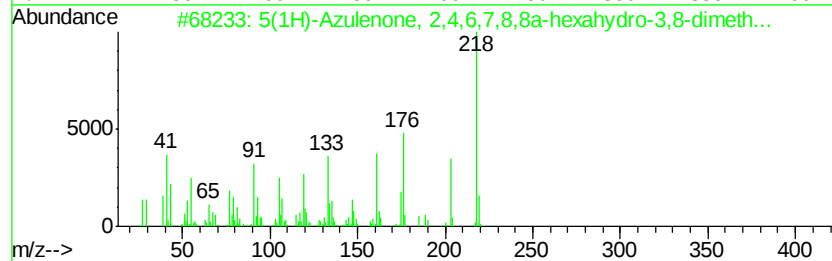
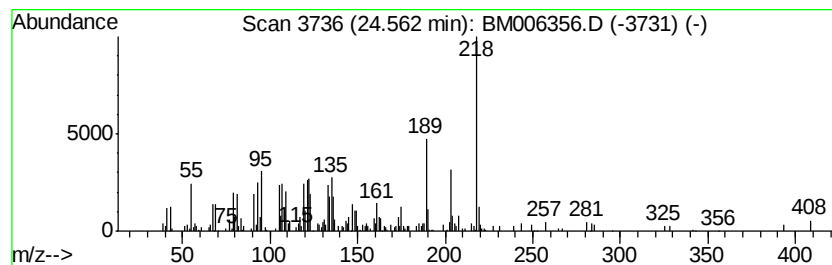
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown-06 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.56	2.79 ng/ul	311982	Perylene-d12	23.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5(1H)-Azulenone, 2,4,6,7,8,8a-he...	218	C15H22O	006754-66-1	49
2		Pyrene, hexadecahydro-	218	C16H26	002435-85-0	38
3		2(1H)Naphthalenone, 3,5,6,7,8,8a...	218	C15H22O	1000188-66-5	38
4		Benzo[b]naphtho[1,2-d]furan	218	C16H10O	000239-30-5	38
5		3-Amino-5-piperonyl-1,2,4-triazole	218	C10H10N4O2	014730-92-8	38



Data Path : V:\HPCHEM1\BNA M\DATA\BM070916\
Data File : BM006356.D
Acq On : 08 Jul 2016 18:42
Operator : UM/SJ
Sample : H3861-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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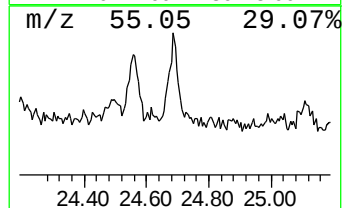
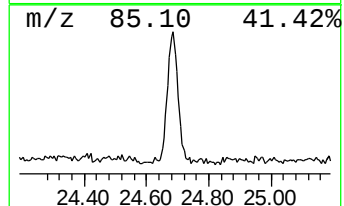
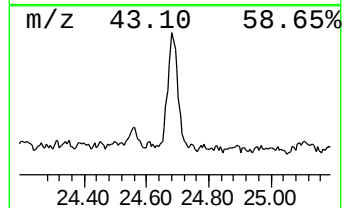
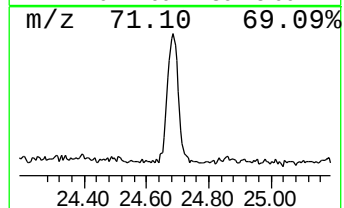
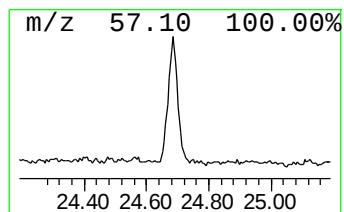
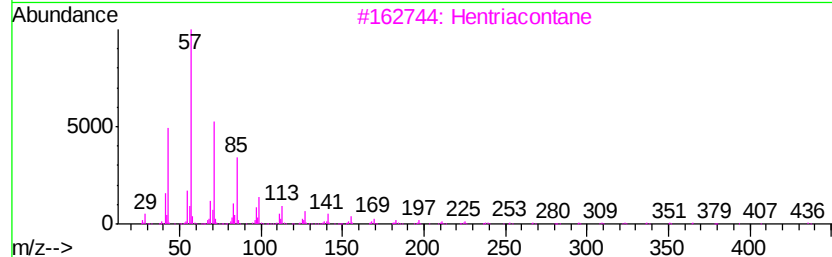
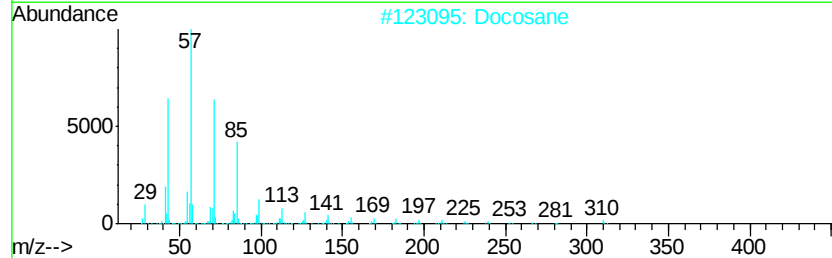
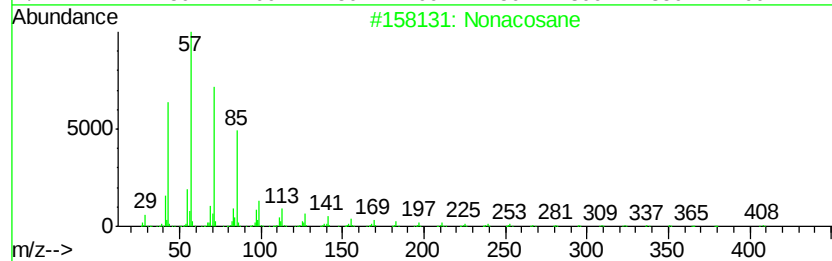
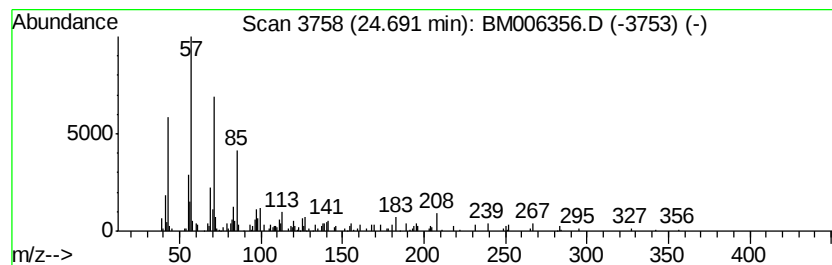
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.69	2.51 ng/ul	280960	Perylene-d12	23.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonacosane	408	C29H60	000630-03-5	86
2			Docosane	310	C22H46	000629-97-0	86
3			Hentriacontane	437	C31H64	000630-04-6	83
4			Tetracosane	338	C24H50	000646-31-1	83
5			Hexacosane	366	C26H54	000630-01-3	83



Data Path : V:\HPCHEM1\BNA_M\DATA\BM070916\
 Data File : BM006356.D
 Acq On : 08 Jul 2016 18:42
 Operator : UM/SJ
 Sample : H3861-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BDJ00

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM070216.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Cyc...	2.83	5.3	ng/ul	312730	1	7.64	1187180	20.0
Bicyclogermacrene	20.37	4.5	ng/ul	537846	5	21.24	2412430	20.0
Azulene, 1,2,3,5,...	20.49	6.2	ng/ul	748851	5	21.24	2412430	20.0
unknown-01	20.56	35.8	ng/ul	4317400	5	21.24	2412430	20.0
Eudesma-4(14),11-...	20.72	12.8	ng/ul	1542000	5	21.24	2412430	20.0
Ethyl 1-thio-.alp...	20.79	33.7	ng/ul	4065240	5	21.24	2412430	20.0
2-Dimethylamino-3...	20.84	30.3	ng/ul	3653860	5	21.24	2412430	20.0
unknown-02	20.90	13.9	ng/ul	1670790	5	21.24	2412430	20.0
unknown-03	21.09	81.1	ng/ul	9777780	5	21.24	2412430	20.0
unknown-04	21.15	123.4	ng/ul	14888000	5	21.24	2412430	20.0
3-Chlorbenz-4-nit...	21.33	10.9	ng/ul	1316900	5	21.24	2412430	20.0
unknown-05	21.45	7.2	ng/ul	867165	5	21.24	2412430	20.0
(DEL) Alkane: Str...	23.96	2.3	ng/ul	257153	6	23.44	2234900	20.0
unknown-06	24.56	2.8	ng/ul	311982	6	23.44	2234900	20.0
(DEL) Alkane: Str...	24.69	2.5	ng/ul	280960	6	23.44	2234900	20.0